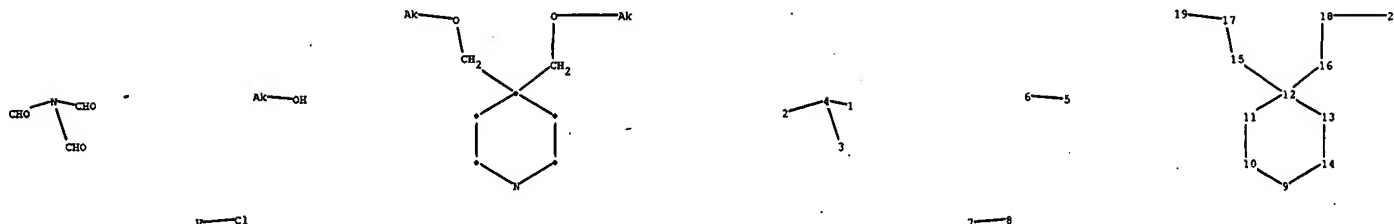


STN Search
 file
 9/19/07



chain nodes :

1 2 3 4 5 6 7 8 15 16 17 18 19 20

ring nodes :

9 10 11 12 13 14

chain bonds :

1-4 2-4 3-4 5-6 7-8 12-15 12-16 15-17 16-18 17-19 18-20

ring bonds :

9-10 9-14 10-11 11-12 12-13 13-14

exact/norm bonds :

1-4 2-4 3-4 5-6 9-10 9-14 10-11 11-12 12-13 13-14 17-19 18-20

exact bonds :

7-8 12-15 12-16 15-17 16-18

Match level :

1:CLASS2:CLASS3:CLASS4:CLASS5:CLASS6:CLASS7:CLASS8:CLASS9:Atom 10:Atom 11:Atom
 12:Atom 13:Atom 14:Atom 15:CLASS16:CLASS17:CLASS18:CLASS19:CLASS20:CLASS

fragments assigned product role:

containing 9

fragments assigned reactant/reagent role:

containing 1

containing 5

containing 7

10/619436 Ketals

=>

Uploading C:\Program Files\Stnexp\Queries\2007 cases\10619436\casreact.str

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 17:14:40 FILE 'CASREACT'

SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED VERIFICATIONS: 0 TO 0

PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1 (0 REACTIONS)

=> s l1 sss full

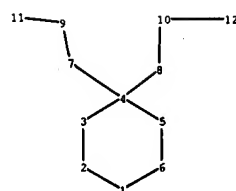
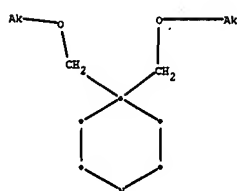
FULL SEARCH INITIATED 17:14:57 FILE 'CASREACT'

SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS

SEARCH TIME: 00:00.01

L3 0 SEA SSS FUL L1 (0 REACTIONS)



chain nodes :

7 8 9 10 11 12

ring nodes :

1 2 3 4 5 6

chain bonds :

4-7 4-8 7-9 8-10 9-11 10-12

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6

exact/norm bonds :

1-2 1-6 2-3 3-4 4-5 5-6 9-11 10-12

exact bonds :

4-7 4-8 7-9 8-10

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS
12:CLASS

10/619436 Ketals

=>

Uploading C:\Program Files\Stnexp\Queries\2007 cases\10619436\triacetonamine
ketal.str

L5 STRUCTURE UPLOADED

=> d 15

L5 HAS NO ANSWERS

L5 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

=> s 15

SAMPLE SEARCH INITIATED 17:19:15 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 52 TO ITERATE

100.0% PROCESSED 52 ITERATIONS

0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**

PROJECTED ITERATIONS: 608 TO 1472

PROJECTED ANSWERS: 0 TO 0

L6 0 SEA SSS SAM L5

=> s 15 sss full

FULL SEARCH INITIATED 17:19:21 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 971 TO ITERATE

100.0% PROCESSED 971 ITERATIONS

0 ANSWERS

SEARCH TIME: 00.00.01

L7 0 SEA SSS FUL L5

10/619436 Ketals

=> d his

(FILE 'HOME' ENTERED AT 17:13:26 ON 19 SEP 2007)

FILE 'CASREACT' ENTERED AT 17:14:03 ON 19 SEP 2007

L1 STRUCTURE UPLOADED

L2 0 S L1

L3 0 S L1 SSS FULL

FILE 'STNGUIDE' ENTERED AT 17:15:51 ON 19 SEP 2007

FILE 'REGISTRY' ENTERED AT 17:17:18 ON 19 SEP 2007

L4 STRUCTURE UPLOADED

FILE 'STNGUIDE' ENTERED AT 17:18:26 ON 19 SEP 2007

FILE 'REGISTRY' ENTERED AT 17:18:46 ON 19 SEP 2007

L5 STRUCTURE UPLOADED

L6 0 S L5

L7 0 S L5 SSS FULL

FILE 'CAPLUS' ENTERED AT 17:20:09 ON 19 SEP 2007

L8 2 S "TRIACETONAMINE KETAL"

FILE 'STNGUIDE' ENTERED AT 17:21:15 ON 19 SEP 2007

L9 0 S "HYDROCHLORIC ACID" (20S) CATALYST

L10 0 S "HYDROGEN CHLORIDE" (20S) CATALYST

L11 0 S "HYDROCHLORIC ACID" (P) CATALYST

L12 0 S "HYDROGEN CHLORIDE" (P) CATALYST

L13 0 S "HYDROGEN CHLORIDE" (P) "TRIACETONAMINE"

L14 0 S "HYDROGEN CHLORIDE" (P) GAS

L15 0 S "HYDROGEN CHLORIDE GAS" OR "GASEOUS HYDROGEN CHLORIDE"

FILE 'HCAPLUS' ENTERED AT 17:30:35 ON 19 SEP 2007

L16 1091 S "HYDROGEN CHLORIDE GAS" OR "GASEOUS HYDROGEN CHLORIDE"

L17 0 S L16 (P) TRIACETONAMINE

L18 1 S "HYDROGEN CHLORIDE" (P) "TRIACETONAMINE"

10/619436 Ketals

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSPTAMJ1621

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

***** Welcome to STN International *****

NEWS 1 JUL 02 Web Page for STN Seminar Schedule - N. America
NEWS 2 JUL 02 LINEDLINE coverage updated
NEWS 3 JUL 02 SCISEARCH enhanced with complete author names
NEWS 4 JUL 02 CHEMCATS accession numbers revised
NEWS 5 JUL 02 CA/CAPLUS enhanced with utility model patents from China
NEWS 6 JUL 16 CAPLUS enhanced with French and German abstracts
NEWS 7 JUL 18 CA/CAPLUS patent coverage enhanced
NEWS 8 JUL 26 USPATFULL/USPAT2 enhanced with IPC reclassification
NEWS 9 JUL 30 USGENE now available on STN
NEWS 10 AUG 06 CAS REGISTRY enhanced with new experimental property tags
NEWS 11 AUG 06 BEILSTEIN updated with new compounds
NEWS 12 AUG 06 FSTA enhanced with new chesaurus edition
NEWS 13 AUG 13 CA/CAPLUS enhanced with additional kind codes for granted patents
NEWS 14 AUG 20 CA/CAPLUS enhanced with CAS indexing in pre-1907 records
NEWS 15 AUG 27 Full-text patent databases enhanced with predefined patent family display formats from INPADOCDB
NEWS 16 AUG 27 USPATOLD now available on STN
NEWS 17 AUG 28 CAS REGISTRY enhanced with additional experimental spectral property data
NEWS 18 SEP 07 STN Anavist, Version 2.0, now available with Derwent World Patents Index
NEWS 19 SEP 13 FORIS renamed to SOPIS
NEWS 20 SEP 13 INPADOCDB enhanced with monthly SDI frequency
NEWS 21 SEP 17 CA/CAPLUS enhanced with printed CA page images from 1967-1998
NEWS 22 SEP 17 CAPLUS coverage extended to include traditional medicine patents

NEWS EXPRESS 19 SEPTEMBER 2007: CURRENT WINDOWS VERSION IS V8.2,
CURRENT MACINTOSH VERSION IS V6.0C(ENG) AND V6.0JC(JP),
AND CURRENT DISCOVER FILE IS DATED 19 SEPTEMBER 2007.

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NEWS LOGIN Welcome Banner and News Items
NEWS IPC8 For general information regarding STN implementation of IPC 8

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Page 1 searched 9/19/07

10/619436 Ketals

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***** STN Columbus *****

FILE 'HOME' ENTERED AT 17:13:26 ON 19 SEP 2007

=> file casreact

COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE ENTRY TOTAL
0.21 0.21

FILE 'CASREACT' ENTERED AT 17:14:03 ON 19 SEP 2007

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FILE CONTENT:1840 - 15 Sep 2007 VOL 147 ISS 13

New CAS Information Use Policies, enter HELP USAGETERMS for details.

* CASREACT now has more than 12 million reactions *

Some CASREACT records are derived from the ZIC/VINITI database (1974-1999) provided by Infochem, INPI data prior to 1986, and Bioinformatics database compiled under the direction of Professor Dr. Klaus Kieslich.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> Uploading C:\Program Files\Stnexp\Queries\2007cases\10619436\casreact.str

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 17:14:40 FILE 'CASREACT'

SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS

SEARCH TIME: 00.00.01

Page 2 searched 9/19/07

10/619436 Ketals

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
PROJECTED VERIFICATIONS: BATCH **COMPLETE**
PROJECTED ANSWERS: 0 TO 0
0 TO 0
L2 0 SEA SSS SAM L1 (0 REACTIONS)
=> s l1 sss full
FULL SEARCH INITIATED 17:14:57 FILE 'CASREACT'
SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS
100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.01
L3 0 SEA SSS FUL L1 (0 REACTIONS)
=> file stng
COST IN U.S. DOLLARS
FULL ESTIMATED COST
FILE 'STNGUIDE' ENTERED AT 17:15:51 ON 19 SEP 2007
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FILE CONTAINS CURRENT INFORMATION.
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=> file reg
COST IN U.S. DOLLARS
FULL ESTIMATED COST
FILE 'REGISTRY' ENTERED AT 17:17:18 ON 19 SEP 2007
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COPYRIGHT (C) 2007 American Chemical Society (ACS)
Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.
STRUCTURE FILE UPDATES: 18 SEP 2007 HIGHEST RN 947490-11-1
DICTIONARY FILE UPDATES: 18 SEP 2007 HIGHEST RN 947490-11-1
New CAS Information Use Policies, enter HELP USAGETERMS for details.
TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007
Please note that search-term pricing does apply when
conducting SmartSELECT searches.
REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:
<http://www.cas.org/support/stngen/stdoc/properties.html>

10/619436 Ketals

=> Uploading C:\Program Files\Stnexp\Queries\2007cases\10619436\triacetonamine
ketals.str
L4 STRUCTURE UPLOADED
=> d l4
L4 HAS NO ANSWERS
L4 STR
L4 *** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
Structure attributes must be viewed using STN Express query preparation.
=> file stng
COST IN U.S. DOLLARS
FULL ESTIMATED COST
FILE 'STNGUIDE' ENTERED AT 17:18:26 ON 19 SEP 2007
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LAST RELOADED: Sep 14, 2007 (20070914/UP).

=> Uploading
THIS COMMAND NOT AVAILABLE IN THE CURRENT FILE
Do you want to switch to the Registry File?
Choice (Y/N):
Switching to the Registry File...
Some commands only work in certain files. For example, the EXPAND
command can only be used to look at the index in a file which has an
index. Enter "HELP COMMANDS" at an arrow prompt (=>) for a list of
commands which can be used in this file.

=> FILE REGISTRY
COST IN U.S. DOLLARS
FULL ESTIMATED COST
FILE 'REGISTRY' ENTERED AT 17:18:46 ON 19 SEP 2007
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provided by InfoChem.
STRUCTURE FILE UPDATES: 18 SEP 2007 HIGHEST RN 947490-11-1
DICTIONARY FILE UPDATES: 18 SEP 2007 HIGHEST RN 947490-11-1
New CAS Information Use Policies, enter HELP USAGETERMS for details.
TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stdoc/properties.html>

```
=>
Uploading C:\Program Files\Stnexp\Queries\2007cases\10619436\triacetonomine
ketal.str
L5 STRUCTURE UPLOADED
```

```
=> d 15
L5 HAS NO ANSWERS
STR
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
```

Structure attributes must be viewed using STN Express query preparation.

```
=> s 15
SAMPLE SEARCH INITIATED 17:19:15 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 52 TO ITERATE
100.0% PROCESSED 52 ITERATIONS
SEARCH TIME: 00.00.01 0 ANSWERS
```

```
FULL FILE PROJECTIONS: ONLINE **COMPLETE**
PROJECTED ITERATIONS: BATCH **COMPLETE**
PROJECTED ANSWERS: 608 TO 1472
0 TO 0
```

```
L6 0 SEA SSS SAM L5
```

```
=> s 15 sss full
FULL SEARCH INITIATED 17:19:21 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 971 TO ITERATE
```

```
100.0% PROCESSED 971 ITERATIONS
SEARCH TIME: 00.00.01 0 ANSWERS
```

```
L7 0 SEA SSS FUL L5
```

```
=> file caplus
COST IN U.S. DOLLARS
FULL ESTIMATED COST
SINCE FILE ENTRY TOTAL
172.55 288.29
```

FILE 'CAPLUS' ENTERED AT 17:20:09 ON 19 SEP 2007
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FILE COVERS 1907 - 19 Sep 2007 VOL 147 ISS 13
FILE LAST UPDATED: 18 Sep 2007 (20070918/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

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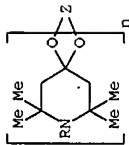
```
=> s "triacetonomine ketal"
260 "TRIACETONAMINE"
2 "TRIACETONAMINES"
261 "TRIACETONAMINE"
("TRIACETONAMINE" OR "TRIACETONAMINES")
10060 "KETAL"
4318 "KETALS"
12133 "KETAL"
```

```
L8 2 "TRIACETONAMINE KETAL"
("KETAL" OR "KETALS")
("TRIACETONAMINE" (W) "KETAL")
```

```
=> d 18 1-2 ibib abs
```

```
L8 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2007 ACS on STN
ACCESSION NUMBER: 1985:504945 CAPLUS
DOCUMENT NUMBER: 103:104945
TITLE: Triacetonomine sugar alcohol ketals
PATENT ASSIGNEE(S): Adeka Argus Chemical Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 5 pp.
CODEN: JKXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:
```

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 60045583	A	19850312	JP 1983-153437	19830823
PRIORITY APPLN. INFO.:			JP 1983-153437	19830823



AB Light-degradable organic materials can be stabilized by title compds. I (R = H, alkyl, alkanoyl; Z = sugar alc. moiety; n = 1, 2). Thus, refluxing 20.0 g xylitol and 33.0 g triacetoneamine-H₂SO₄ gave a monoketal (I; R = H, Z = xylitol residue, n = 1) (II). Press-molded sheets prepared from polypropylene 100, stearyl μ -(3,5-di-tert-butyl-4-hydroxyphenyl)propionate 0.2, and II 0.3 parts had Hg lamp stability 640 h, vs. 150 h using 8-aza-7,9,9-tetramethyl-1,4-dioxaspiro[4.5]decane.

L8 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2007 ACS ON STN

ACCESSION NUMBER: 1974:27899 CAPLUS

DOCUMENT NUMBER: 80:27899

TITLE: Triacetoneamine ketal stabilizers

INVENTOR(S): Murayama, Keisuke; Toda, Toshimasa; Mori, Eiko; Matsui, Katsuki; Kurumada, Tomoyuki; Ohta, Noriyuki; Watanabe, Ichiro

PATENT ASSIGNEE(S): Sankyo Co., Ltd.

SOURCE: Ger. Offen., 20 pp.

CODEN: GWXXBX

DOCUMENT TYPE: Patent

LANGUAGE: German

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
DE 2203533	A1	19730816	DE 1972-2203533	19720121
PRIORITY APPLN. INFO.: DE 1972-2203533 A 19720121				
AB The ketals I (R = R1 = Bu or n-C8H17, R2 = o-phenylene (II) or CH2CH2) were prepared by ketalization of triacetoneamine (III) and used as heat and light stabilizers in polymers, e.g. polypropylene (IV) [9003-07-0], nylon 6 [25038-54-4], or polyurethanes. Thus, refluxing III and o-(HO)2C6H4 in C6H6 containing p-MeC6H4SO3H gave 2,2,6,6-tetramethyl-4,4-(o-phenylenedioxy)piperidine (II) [36793-29-0]. Samples from 100 parts IV and 0.25 part II turned brittle on heating at 45 deg. under uv irradiation) after 1000 hr vs. 100 hr for IV containing no II.				

=> fill stng

COST IN U.S. DOLLARS

FULL ESTIMATED COST	SINCE FILE		TOTAL SESSION
	ENTRY	10.62	
298.91			

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE		TOTAL SESSION
	ENTRY	-1.56	

CA SUBSCRIBER PRICE

TOTAL SESSION	TOTAL SESSION

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FILE CONTAINS CURRENT INFORMATION.

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=> d his

(FILE 'HOME' ENTERED AT 17:13:26 ON 19 SEP 2007)

FILE 'CASREACT' ENTERED AT 17:14:03 ON 19 SEP 2007

STRUCTURE UPLOADED

L1 0 S L1

L2 0 S L1 SSS FULL

FILE 'STNGUIDE' ENTERED AT 17:15:51 ON 19 SEP 2007

FILE 'REGISTRY' ENTERED AT 17:17:18 ON 19 SEP 2007

STRUCTURE UPLOADED

FILE 'STNGUIDE' ENTERED AT 17:18:26 ON 19 SEP 2007

FILE 'REGISTRY' ENTERED AT 17:18:46 ON 19 SEP 2007

STRUCTURE UPLOADED

L5 0 S L5

L6 0 S L5 SSS FULL

FILE 'CAPLUS' ENTERED AT 17:20:09 ON 19 SEP 2007

2 S "TRIACETONAMINE KETAL"

FILE 'STNGUIDE' ENTERED AT 17:21:15 ON 19 SEP 2007

=> s "hydrochloric acid" (20s) catalyst

0 "HYDROCHLORIC"

7 "ACID"

1 "ACIDS"

7 "ACID"

("ACID" OR "ACIDS")

0 "HYDROCHLORIC ACID"

("HYDROCHLORIC" (W) "ACID")

7 CATALYST

6 CATALYSTS

8 CATALYST

(CATALYST OR CATALYSTS)

FILE 'STNGUIDE' ENTERED AT 17:21:15 ON 19 SEP 2007

0 "HYDROCHLORIC ACID" (20S) CATALYST

=> s "hydrogen chloride" (20s) catalyst

4 "HYDROGEN"

0 "CHLORIDE"

0 "HYDROGEN CHLORIDE"

("HYDROGEN" (W) "CHLORIDE")

7 CATALYST

6 CATALYSTS

8 CATALYST

(CATALYST OR CATALYSTS)

FILE 'STNGUIDE' ENTERED AT 17:21:15 ON 19 SEP 2007

0 "HYDROGEN CHLORIDE" (20S) CATALYST

=> s "hydrochloric acid" (p) catalyst

0 "HYDROCHLORIC"

10/619436 Ketals

```

7 "ACID"
1 "ACIDS"
7 "ACID"
  ("ACID" OR "ACIDS")
0 "HYDROCHLORIC ACID"
  ("HYDROCHLORIC" (W) "ACID")
7 CATALYST
6 CATALYSTS
8 CATALYST
  (CATALYST OR CATALYSTS)
L11 0 "HYDROCHLORIC ACID" (P) CATALYST
=> s "hydrogen chloride" (p) catalyst
4 "HYDROGEN"
0 "CHLORIDE"
0 "HYDROGEN CHLORIDE"
  ("HYDROGEN" (W) "CHLORIDE")
7 CATALYST
6 CATALYSTS
8 CATALYST
  (CATALYST OR CATALYSTS)
L12 0 "HYDROGEN CHLORIDE" (P) CATALYST
=> s "hydrogen chloride" (p) "triacetanamine"
4 "HYDROGEN"
0 "CHLORIDE"
0 "HYDROGEN CHLORIDE"
  ("HYDROGEN" (W) "CHLORIDE")
0 "TRIACETANAMINE"
0 "HYDROGEN CHLORIDE" (P) "TRIACETANAMINE"
L13
=> s "hydrogen chloride" (p) gas
4 "HYDROGEN"
0 "CHLORIDE"
0 "HYDROGEN CHLORIDE"
  ("HYDROGEN" (W) "CHLORIDE")
8 GAS
3 GASES
10 GAS
  (GAS OR GASES)
L14 0 "HYDROGEN CHLORIDE" (P) GAS
=> s "hydrogen chloride gas" or "gaseous hydrogen chloride"
4 "HYDROGEN"
0 "CHLORIDE"
8 "GAS"
3 "GASES"
10 "GAS"
  ("GAS" OR "GASES")
0 "HYDROGEN CHLORIDE GAS"
  ("HYDROGEN" (W) "CHLORIDE" (W) "GAS")
0 "GASEOUS"
4 "HYDROGEN"
0 "CHLORIDE"
0 "GASEOUS HYDROGEN CHLORIDE"
  ("GASEOUS" (W) "HYDROGEN" (W) "CHLORIDE")
L15 0 "HYDROGEN CHLORIDE GAS" OR "GASEOUS HYDROGEN CHLORIDE"

```

Page 9 searched 9/19/07

10/619436 Ketals

```

=> d his
(FILE 'HOME' ENTERED AT 17:13:26 ON 19 SEP 2007)
FILE 'CASREACT' ENTERED AT 17:14:03 ON 19 SEP 2007
STRUCTURE UPLOADED
0 S L1
0 S L1 SSS FULL

FILE 'STNGUIDE' ENTERED AT 17:15:51 ON 19 SEP 2007
FILE 'REGISTRY' ENTERED AT 17:17:18 ON 19 SEP 2007
STRUCTURE UPLOADED

FILE 'STNGUIDE' ENTERED AT 17:18:26 ON 19 SEP 2007
FILE 'REGISTRY' ENTERED AT 17:18:46 ON 19 SEP 2007
STRUCTURE UPLOADED
0 S L5
0 S L5 SSS FULL

FILE 'CAPLUS' ENTERED AT 17:20:09 ON 19 SEP 2007
2 S "TRIACETAMINE KETAL"

FILE 'STNGUIDE' ENTERED AT 17:21:15 ON 19 SEP 2007
0 S "HYDROCHLORIC ACID" (20S) CATALYST
0 S "HYDROGEN CHLORIDE" (20S) CATALYST
0 S "HYDROCHLORIC ACID" (P) CATALYST
0 S "HYDROGEN CHLORIDE" (P) CATALYST
0 S "HYDROGEN CHLORIDE" (P) CATALYST
0 S "HYDROGEN CHLORIDE" (P) GAS
0 S "HYDROGEN CHLORIDE" (P) GAS
0 S "HYDROGEN CHLORIDE GAS" OR "GASEOUS HYDROGEN CHLORIDE"

=> file hcapl
COST IN U.S. DOLLARS
FULL ESTIMATED COST
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)
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Page 10 searched 9/19/07

FILE COVERS 1907 - 19 Sep 2007 VOL 147 ISS 13
FILE LAST UPDATED: 18 Sep 2007 (20070918/ED)

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s "hydrogen chloride gas" or "gaseous hydrogen chloride"

1015422 "HYDROGEN"
6041 "HYDROGENS"
1018782 "HYDROGEN" OR "HYDROGENS"
1161611 "CHLORIDE"
159679 "CHLORIDES"
1235117 "CHLORIDE"
1609761 "GAS"
534142 "GASES"
1797109 "GAS"
716 "HYDROGEN CHLORIDE GAS"
174183 "GASEOUS"
1015422 "HYDROGEN"
6041 "HYDROGENS"
1018782 "HYDROGEN" OR "HYDROGENS"
1161611 "CHLORIDE"
159679 "CHLORIDES"
1235117 "CHLORIDE"
407 "GASEOUS HYDROGEN CHLORIDE"
1091 "HYDROGEN CHLORIDE GAS" OR "GASEOUS HYDROGEN CHLORIDE"

L16 1091 "HYDROGEN CHLORIDE GAS" OR "GASEOUS HYDROGEN CHLORIDE"

=> l16 (P) triacetanamine

L16 IS NOT A RECOGNIZED COMMAND

The previous command name entered was not recognized by the system.

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=> s l16 (P) triacetanamine
260 TRIACETONAMINE
261 TRIACETONAMINES
0 L16 (P) TRIACETONAMINE OR TRIACETONAMINES

L17

=> s "hydrogen chloride" (P) "triacetanamine"
1015422 "HYDROGEN"
6041 "HYDROGENS"
1018782 "HYDROGEN" OR "HYDROGENS"
1161611 "CHLORIDE"
159679 "CHLORIDES"
1235117 "CHLORIDE" OR "CHLORIDES"

33237 "HYDROGEN CHLORIDE"
("HYDROGEN" (W) "CHLORIDE")

260 "TRIACETONAMINE"

2 "TRIACETONAMINES"

261 "TRIACETONAMINE"

("TRIACETONAMINE" OR "TRIACETONAMINES")

1 "HYDROGEN CHLORIDE" (P) "TRIACETONAMINE"

=> d l18 ibib abs

L18 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1906:88385 HCAPLUS

DOCUMENT NUMBER: 0:88385

TITLE: The oximes of the cyclic acetone bases:

paramidotrimeethylpiperidine

Harties, Carl D.

SOURCE: Berichte der Deutschen Chemischen Gesellschaft (1896),

29, 521-9

From: J. Chem. Soc., Abstr. 70, 1, 317-8 1896

CODEN: BDCGAS

DOCUMENT TYPE:

LANGUAGE: Unavailable

AB Triacetanaminoxime, C₉H₁₈N₂O, is obtained by the action of hydroxylamine

on triacetanamin, and crystallises in large, white, six-sided

prisms, melting at 152-153°; it forms crystalline salts with

hydrochloric and sulphuric acids. Benzylidenediacetonaminoxime,

C₁₃H₁₈N₂O, crystallises in lustrous, four-sided tablets, melting at

140-141°, and is only sparingly soluble in boiling water.

Vinylidiacetonaminoxime, (the original abstract includes an equation),

crystallises in translucent, four-sided tablets, melting at

150-151°. When reduced with alcoholic hydrogen

chloride, zinc dust, and a little water, it is converted into

4-amido-2:2:6-trimethylpiperidine, which forms a crystalline mass, melts

at 25-26°, and boils at 60° (pressure = 7.5 mm.). The base

has a faint odour of piperidine, and rapidly combines with the carbonic

anhydride of the air, forming a carbamate. The hydriodide of the base

crystallises in fascicular groups of white prisms, and is moderately

soluble in water; the hydrochloride is readily soluble in water; the

autochloride crystallises in red, oblique, six-sided tablets, and is

sparingly soluble in water, whilst the platinochloride and picrate are

also crystalline. The base forms both a normal and an acid oxalate, the

latter being very hygroscopic. 4-acetamido-2:2:6-trimethylpiperidine

crystallises in cubes, melting at 206-207°; it is strongly basic,

and forms an autochloride, melting at 235° with decomposition. The

diacetyl compound, (the original abstract includes an equation), is formed

when the base is heated with excess of acetic anhydride at 160°; it

forms small prisms, melts at 88-89°, boils at 160-170°

(pressure = 8 mm.), and has basic properties, forming a crystalline

autochloride. This diacetyl compound is accompanied by another basic

substance, which boils at about 200° (pressure = 8 mm.), and is

probably an anhydro-derivative. The base does not yield a diazo-compound

with sodium nitrite and an acid, whilst with amyllic nitrite it yields a

nitroso-derivative, the imido-group having taken part in the reaction.

When heated with chloroform and alcoholic potash, no carbonylamine

derivative is produced. The base reacts with carbon bisulphide, forming a

thiocarbamate, C₉H₁₈N₂S₂, which crystallises from water in prisms and

melts at 187-188°. When this salt is treated with mercuric

chloride, it yields the hydrochloride of a new base. The latter

crystallises in small prisms and melts at 79-80°. It has not the smell or other properties of a thiocarbimide, and probably has the constitution (the original abstract includes an equation).

=> d his

(FILE 'HOME' ENTERED AT 17:13:26 ON 19 SEP 2007)

FILE 'CASREACT' ENTERED AT 17:14:03 ON 19 SEP 2007
STRUCTURE UPLOADED

L1 0 S L1
L2 0 S L1
L3 0 S L1 SSS FULL

FILE 'STNGUIDE' ENTERED AT 17:15:51 ON 19 SEP 2007

FILE 'REGISTRY' ENTERED AT 17:17:18 ON 19 SEP 2007
STRUCTURE UPLOADED

L4

FILE 'STNGUIDE' ENTERED AT 17:18:26 ON 19 SEP 2007

FILE 'REGISTRY' ENTERED AT 17:18:46 ON 19 SEP 2007
STRUCTURE UPLOADED

L5 0 S L5
L6 0 S L5
L7 0 S L5 SSS FULL

FILE 'CAPLUS' ENTERED AT 17:20:09 ON 19 SEP 2007
2 S "TRIACETONAMINE RETAL"

L8

FILE 'STNGUIDE' ENTERED AT 17:21:15 ON 19 SEP 2007
0 S "HYDROCHLORIC ACID" (20S) CATALYST

L9 0 S "HYDROGEN CHLORIDE" (20S) CATALYST
L10 0 S "HYDROCHLORIC ACID" (20S) CATALYST
L11 0 S "HYDROCHLORIC ACID" (P) CATALYST
L12 0 S "HYDROGEN CHLORIDE" (P) CATALYST
L13 0 S "HYDROGEN CHLORIDE" (P) "TRIACETONAMINE"
L14 0 S "HYDROGEN CHLORIDE" (P) GAS
L15 0 S "HYDROGEN CHLORIDE GAS" OR "GASEOUS HYDROGEN CHLORIDE"

FILE 'HCAPLUS' ENTERED AT 17:30:35 ON 19 SEP 2007
1091 S "HYDROGEN CHLORIDE GAS" OR "GASEOUS HYDROGEN CHLORIDE"

L16 0 S L16 (P) TRIACETONAMINE
L17 1 S "HYDROGEN CHLORIDE" (P) "TRIACETONAMINE"

L18

=> fil slng

COST IN U.S. DOLLARS

SINCE FILE ENTRY	TOTAL SESSION
15.83	315.70

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE ENTRY	TOTAL SESSION
-0.78	-2.34

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LAST RELOADED: Sep 14, 2007 (20070914/UP).

EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	271	"564/502" CCL5	US-PGPUB; USPAT; USOCR	OR	ON	2007/09/19 16:18
L2	513	"544/406" CCL5	US-PGPUB; USPAT; USOCR	OR	ON	2007/09/19 16:18
L3	794	L1 same20 L2	US-PGPUB; USPAT; USOCR	OR	ON	2007/09/19 16:18
L4	125	tracetbonamine	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:19
L5	12	ketal5 same L4	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:19
L6	125	tracetbonamine	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:37
L7	2	("3959298") URPN	USPAT	OR	ON	2007/09/19 16:23
L8	1	("4734502") URPN	USPAT	OR	ON	2007/09/19 16:23
L9	0	("4831146") URPN	USPAT	OR	ON	2007/09/19 16:23
L10	7	tracetbonam3 near3 ketal	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:34
L11	201185	((hydrogen chloride) or (hydrochloric acid)) near3 (gas or air or vapor)	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:36
L12	201301	L11 same20 L4	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:36
L13	0	L11 same20 L4	US-PGPUB; USPAT; EPO; JPO; DERWENT	AND	ON	2007/09/19 16:36
L14	2	L11 near20 L4	US-PGPUB; USPAT; EPO; JPO; DERWENT	AND	ON	2007/09/19 16:36

EAST Search History

L15	0	tracetbonamine.dm. and ketal.dm.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:40
L16	0	tracetbonamine.dm. and ketal.dm.	US-PGPUB; USPAT; EPO; JPO; DERWENT	ADJ	ON	2007/09/19 16:39
L17	156	L11 and ketal.cim.	US-PGPUB; USPAT; EPO; JPO; DERWENT	ADJ	ON	2007/09/19 16:39
L18	0	L17 and L6	US-PGPUB; USPAT; EPO; JPO; DERWENT	ADJ	ON	2007/09/19 16:39
L19	2353	tracetbonamine.dm. or ketal.cim.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:40
L20	158	L11 and L19	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:48
L21	189	ketal510 near10 (tracetbonamine or piperidone)	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:49
L22	189	(ketal510) near10 (tracetbonamine or piperidone)	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:57
L23	186	(ketal510) near10 piperidone	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:53
L24	8	(ketal510) near10 tracetbonamine	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/09/19 16:57
S1	3	("3963730") or ("4250312") or ("3940401") PN.	US-PGPUB; USPAT	OR	OFF	2006/12/05 20:29
S2	269	"564/502" CCL5	US-PGPUB; USPAT; USOCR	OR	ON	2007/09/19 16:18
S3	459	"544/406" CCL5	US-PGPUB; USPAT; USOCR	OR	ON	2006/12/05 18:00

EAST Search History

S4	25	(OLIVER) near2 (MEYER)). INV.	US-PGPUB; USPAT; USOCR	OR	ON	2006/12/05 18:00
S5	26	((OLIVER) near2 (MEYER)). INV.	EPO; JPO; DERWENT	OR	ON	2006/12/05 18:00
S6	1	((RENATE) near2 (UHLENBERG)). INV.	US-PGPUB; USPAT; USOCR	OR	ON	2006/12/05 18:02
S7	0	((RENATE) near2 (UHLENBERG)). INV.	EPO; JPO; DERWENT	OR	ON	2006/12/05 18:02
S8	4	((MICHAEL) near2 (KORELL)). INV.	US-PGPUB; USPAT; USOCR	OR	ON	2006/12/05 18:02
S9	7	((MICHAEL) near2 (KORELL)). INV.	EPO; JPO; DERWENT	OR	ON	2006/12/05 18:03
S10	4	(*4250312*). URPN.	USPAT	OR	ON	2006/12/05 18:23
S11	0	(*4900833*). URPN.	USPAT	OR	ON	2006/12/05 18:28
S12	4	(*4250312*). URPN.	USPAT	OR	ON	2006/12/05 20:28
S13	125	tracetamine	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:55
S15	12	ketals same S13	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:30
S16	791	hydroxyl adj. derivative	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:49
S17	2	S13 same S16	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:38
S18	24509	ketals	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:39
S19	12	S13 same S18	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:40
S20	4	glycerol near5 acid adj. catalysis	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:45

EAST Search History

S21	717	glycerol near5 ketone	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:45
S22	29	S21 same catalys	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:46
S23	27	S16 same ketone	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:50
S24	9	S16 near5 ketone	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:50
S25	59471	hydrogen adj. chloride	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:55
S26	7	S13 and S18 and S25	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:57
S27	24	S13 and S18	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 21:10
S28	2933	S18 and S25	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:58
S29	7	S27 and S28	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 20:58
S30	45181	hydrocarbon adj. solvs	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 21:11
S31	10587	alkalys adj. metal adj. alkoxids	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 21:12
S32	1317	alkalys adj. earth adj. metal adj. alkoxids	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 21:12

EAST Search History

S33	0	S13 same S30	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 21:21
S34	0	S13 near S30	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 21:21
S35	86	S18 same S30	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2006/12/05 21:22
S36	41	("3862100");URPN	USPAT	OR	ON	2006/12/05 21:43
S39	1	JP-48055938-\$.dtd.	US-PGPUB; USPAT; EPO; JPO; DERWENT	OR	ON	2007/01/30 12:11
S40	1	("3839273");PN	US-PGPUB; USPAT	OR	OFF	2007/01/30 12:11
S41	9	("3839273");URPN	USPAT	OR	ON	2007/01/30 12:14